Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 07:24 PM GMT

PDB ID	:	4U4A
Title	:	Complex Structure of BRCA1 BRCT with singly phospho Abraxas
Authors	:	Badgujar, D.; Hosur, M.V.; Varma, A.K.
Deposited on	:	2014-07-23
Resolution	:	3.51 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see references ⓘ) were used in the production of this report:

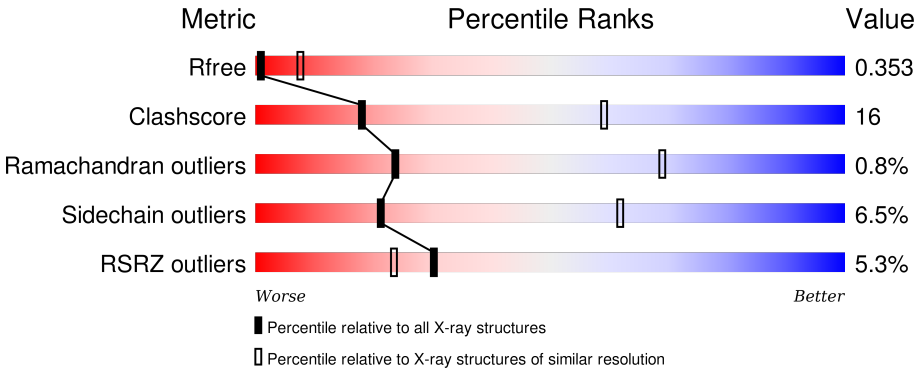
MolProbity	:	4.02b-467
Mogul	:	1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026982
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20026982

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.51 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	1089 (3.64-3.40)
Clashscore	102246	1197 (3.64-3.40)
Ramachandran outliers	100387	1159 (3.64-3.40)
Sidechain outliers	100360	1160 (3.64-3.40)
RSRZ outliers	91569	1096 (3.64-3.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	214	70% 27% ·
1	B	214	6% 71% 26% ·
1	C	214	10% 70% 27% ··
2	D	11	18% 27% 9% 27% 18%
2	E	11	9% 45% 36% 18%

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Mol	Chain	Length	Quality of chain
2	F	11	

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 5332 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Breast cancer type 1 susceptibility protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	214	Total	C	N	O	S	0	0	0
			1709	1093	293	309	14			
1	B	214	Total	C	N	O	S	0	0	0
			1712	1094	294	310	14			
1	C	212	Total	C	N	O	S	0	0	0
			1694	1084	290	306	14			

- Molecule 2 is a protein called BRCA1-A complex subunit Abraxas.

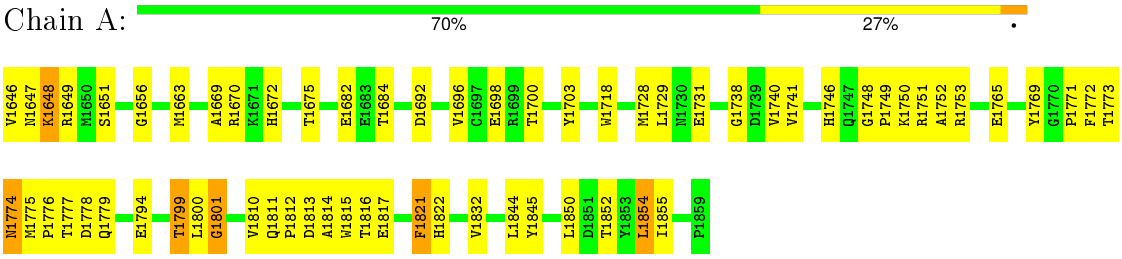
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	9	Total	C	N	O	P	0	0	0
			69	40	12	16	1			
2	E	11	Total	C	N	O	P	0	0	0
			91	57	14	19	1			
2	F	7	Total	C	N	O	P	0	0	0
			57	36	7	13	1			

3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

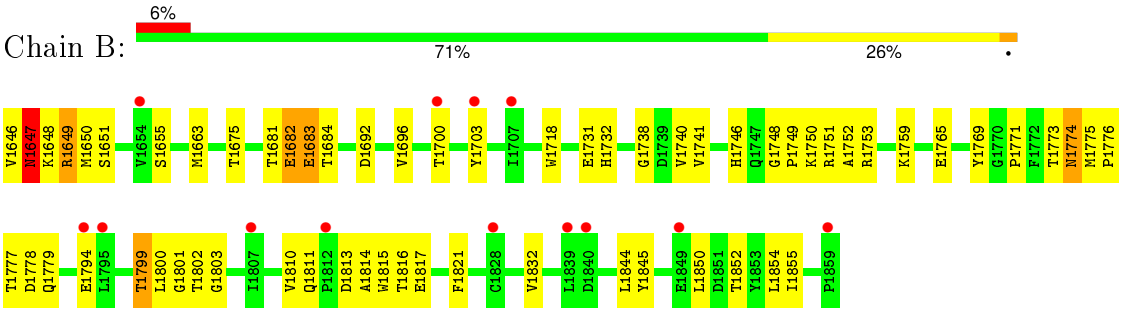
- Molecule 1: Breast cancer type 1 susceptibility protein

Chain A:



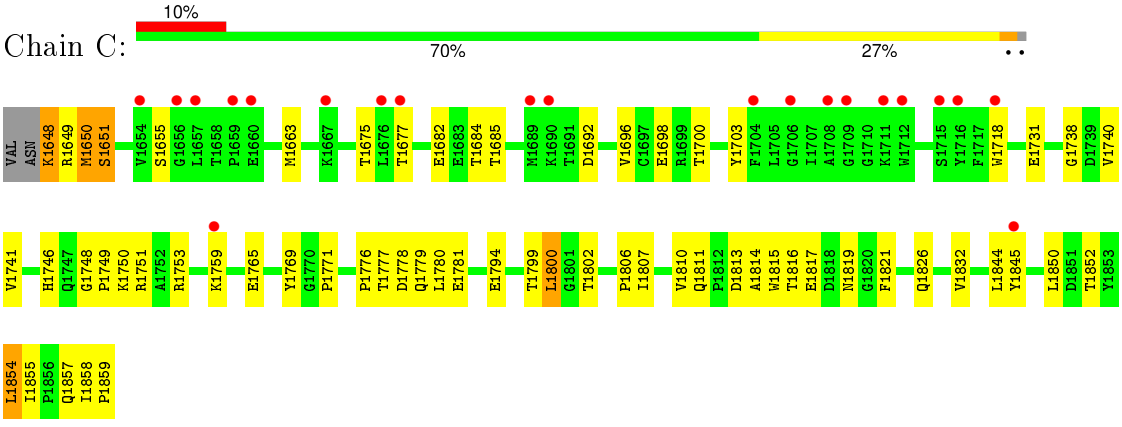
- Molecule 1: Breast cancer type 1 susceptibility protein

Chain B:



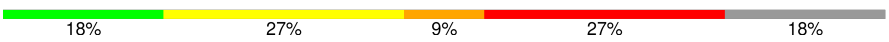
- Molecule 1: Breast cancer type 1 susceptibility protein

Chain C:

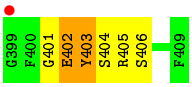


- Molecule 2: BRCA1-A complex subunit Abraxas

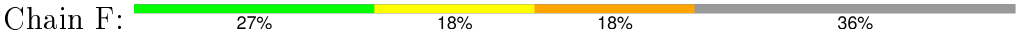
Chain D:

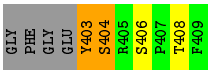


• Molecule 2: BRCA1-A complex subunit Abraxas



• Molecule 2: BRCA1-A complex subunit Abraxas





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α, β, γ	187.50Å 187.50Å 85.35Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	59.89 – 3.51 59.82 – 3.51	Depositor EDS
% Data completeness (in resolution range)	75.6 (59.89-3.51) 75.7 (59.82-3.51)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.56 (at 3.49Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.301 , 0.362 0.295 , 0.353	Depositor DCC
R_{free} test set	754 reflections (5.37%)	DCC
Wilson B-factor (Å^2)	155.6	Xtriage
Anisotropy	0.593	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 129.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Outliers	0 of 14795 reflections	Xtriage
F_o,F_c correlation	0.88	EDS
Total number of atoms	5332	wwPDB-VP
Average B, all atoms (Å^2)	166.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.02% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SEP

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\#\|Z\| > 5$	RMSZ	$\#\|Z\| > 5$
1	A	0.61	0/1750	0.77	0/2374
1	B	0.53	0/1753	0.74	1/2378 (0.0%)
1	C	0.47	0/1735	0.69	0/2353
2	D	1.01	0/59	1.56	2/75 (2.7%)
2	E	0.60	0/83	0.70	0/107
2	F	0.91	0/48	1.21	1/62 (1.6%)
All	All	0.55	0/5428	0.75	4/7349 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	2
1	C	0	1
All	All	0	4

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	D	403	TYR	CB-CA-C	7.05	124.51	110.40
2	F	403	TYR	CA-C-N	5.92	130.23	117.20
2	D	404	SER	N-CA-C	5.77	126.57	111.00
1	B	1647	ASN	N-CA-C	5.33	125.38	111.00

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1801	GLY	Peptide
1	B	1646	VAL	Peptide
1	B	1647	ASN	Peptide
1	C	1650	MET	Peptide

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1709	0	1689	54	1
1	B	1712	0	1692	46	0
1	C	1694	0	1674	58	0
2	D	69	0	55	8	0
2	E	91	0	73	14	1
2	F	57	0	42	3	0
All	All	5332	0	5225	166	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 16.

All (166) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1651:SER:OG	1:B:1684:THR:HA	1.24	1.36
1:C:1776:PRO:CG	1:C:1779:GLN:HG2	1.60	1.29
1:B:1765:GLU:OE2	1:B:1799:THR:OG1	1.55	1.23
1:A:1765:GLU:OE2	1:A:1799:THR:OG1	1.59	1.17
1:B:1648:LYS:HB2	1:B:1649:ARG:HA	1.21	1.09
1:B:1655:SER:OG	2:E:406:SEP:O2P	1.69	1.09
1:C:1776:PRO:HG2	1:C:1779:GLN:CG	1.84	1.06
1:B:1765:GLU:CD	1:B:1799:THR:OG1	1.98	1.01
1:B:1651:SER:OG	1:B:1684:THR:CA	2.11	0.97
1:A:1765:GLU:CD	1:A:1799:THR:OG1	2.03	0.95
1:C:1765:GLU:OE2	1:C:1799:THR:OG1	1.84	0.95
1:C:1776:PRO:CG	1:C:1779:GLN:CG	2.46	0.91
2:D:401:GLY:O	2:D:403:TYR:N	2.03	0.91
1:C:1776:PRO:HG2	1:C:1779:GLN:HG2	0.90	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1648:LYS:HG3	1:C:1854:LEU:HD22	1.56	0.86
1:B:1648:LYS:HB2	1:B:1649:ARG:CA	2.07	0.84
1:B:1648:LYS:HG3	1:B:1649:ARG:HB3	1.58	0.83
1:B:1648:LYS:CB	1:B:1649:ARG:HA	2.07	0.83
1:A:1773:THR:CB	1:A:1774:ASN:OD1	2.32	0.78
1:A:1778:ASP:OD1	1:A:1779:GLN:N	2.17	0.77
1:C:1778:ASP:OD1	1:C:1779:GLN:N	2.16	0.77
2:E:405:ARG:HG3	2:E:406:SEP:N	2.02	0.74
2:D:404:SER:O	2:D:405:ARG:HB2	1.86	0.74
1:A:1648:LYS:HE2	1:C:1857:GLN:HG3	1.70	0.74
1:B:1778:ASP:OD1	1:B:1779:GLN:N	2.20	0.74
1:A:1776:PRO:HG2	1:A:1779:GLN:HG2	1.70	0.74
1:A:1647:ASN:HB2	1:A:1648:LYS:HD3	1.69	0.73
2:E:402:GLU:HB3	2:E:403:TYR:CE2	2.27	0.69
1:A:1773:THR:OG1	1:A:1774:ASN:OD1	2.11	0.69
1:B:1801:GLY:O	1:B:1803:GLY:N	2.26	0.68
1:A:1854:LEU:HD22	1:B:1647:ASN:O	1.94	0.67
1:B:1765:GLU:OE1	1:B:1799:THR:OG1	2.11	0.67
2:F:403:TYR:N	2:F:404:SER:HB2	2.10	0.67
1:A:1765:GLU:OE1	1:A:1799:THR:OG1	2.13	0.67
1:A:1773:THR:HB	1:A:1774:ASN:OD1	1.96	0.66
2:E:402:GLU:HB3	2:E:403:TYR:CZ	2.31	0.65
1:C:1649:ARG:O	1:C:1650:MET:CE	2.45	0.65
1:C:1800:LEU:N	1:C:1800:LEU:HD23	2.12	0.65
1:A:1771:PRO:HD2	1:A:1815:TRP:NE1	2.11	0.64
1:C:1655:SER:OG	2:F:406:SEP:O2P	2.15	0.63
2:E:405:ARG:CG	2:E:406:SEP:N	2.62	0.63
1:A:1648:LYS:CE	1:C:1857:GLN:HG3	2.29	0.62
1:C:1776:PRO:HB2	1:C:1778:ASP:OD1	2.00	0.62
1:B:1648:LYS:CG	1:B:1649:ARG:HB3	2.30	0.61
1:C:1776:PRO:HD2	1:C:1779:GLN:HB2	1.82	0.61
1:A:1774:ASN:OD1	1:A:1774:ASN:N	2.34	0.60
1:C:1771:PRO:HD2	1:C:1815:TRP:NE1	2.17	0.58
1:A:1651:SER:HA	1:A:1675:THR:HB	1.85	0.58
1:B:1682:GLU:O	1:B:1684:THR:N	2.36	0.57
1:B:1776:PRO:HB2	1:B:1778:ASP:OD1	2.04	0.57
1:C:1650:MET:O	1:C:1685:THR:OG1	2.20	0.56
1:A:1773:THR:C	1:A:1774:ASN:OD1	2.44	0.56
1:C:1776:PRO:CD	1:C:1779:GLN:CG	2.84	0.56
1:A:1776:PRO:HG2	1:A:1778:ASP:OD1	2.06	0.56
1:B:1771:PRO:HD2	1:B:1815:TRP:NE1	2.20	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1648:LYS:CB	1:B:1649:ARG:CA	2.76	0.56
1:B:1651:SER:HA	1:B:1675:THR:HB	1.89	0.55
2:E:402:GLU:CB	2:E:403:TYR:CE2	2.89	0.55
1:A:1800:LEU:HD23	1:A:1801:GLY:N	2.22	0.55
1:C:1765:GLU:CD	1:C:1799:THR:OG1	2.45	0.55
1:C:1811:GLN:O	1:C:1814:ALA:HB3	2.07	0.55
1:B:1682:GLU:CD	1:B:1682:GLU:H	2.10	0.54
1:A:1647:ASN:HB2	1:A:1648:LYS:CD	2.36	0.54
1:C:1777:THR:O	1:C:1780:LEU:HB3	2.08	0.53
1:A:1811:GLN:O	1:A:1814:ALA:HB3	2.08	0.53
2:E:403:TYR:HB3	2:E:404:SER:HA	1.90	0.53
1:A:1648:LYS:HE2	1:C:1857:GLN:CG	2.37	0.52
2:D:401:GLY:C	2:D:403:TYR:N	2.63	0.52
1:C:1651:SER:O	1:C:1684:THR:HA	2.10	0.52
1:B:1811:GLN:O	1:B:1814:ALA:HB3	2.09	0.52
1:A:1648:LYS:N	1:A:1648:LYS:HD3	2.25	0.51
1:C:1650:MET:HB3	1:C:1685:THR:HG23	1.93	0.51
1:C:1813:ASP:N	1:C:1813:ASP:OD1	2.43	0.51
2:E:402:GLU:C	2:E:403:TYR:CG	2.85	0.51
1:B:1682:GLU:O	1:B:1683:GLU:C	2.49	0.50
1:A:1812:PRO:HB2	1:B:1732:HIS:NE2	2.26	0.50
1:C:1778:ASP:HA	1:C:1781:GLU:HB2	1.93	0.50
2:E:402:GLU:HB3	2:E:403:TYR:CD2	2.46	0.50
2:D:402:GLU:O	2:D:403:TYR:C	2.48	0.50
2:E:403:TYR:CB	2:E:404:SER:HA	2.41	0.49
1:C:1698:GLU:OE2	2:F:408:THR:HG21	2.11	0.49
1:A:1813:ASP:N	1:A:1813:ASP:OD1	2.44	0.49
1:B:1813:ASP:OD1	1:B:1813:ASP:N	2.44	0.49
1:B:1681:THR:O	1:B:1682:GLU:C	2.51	0.48
1:B:1692:ASP:OD1	1:B:1696:VAL:N	2.42	0.48
2:D:403:TYR:CB	2:D:404:SER:CB	2.91	0.48
1:A:1740:VAL:HG23	1:A:1741:VAL:HG23	1.95	0.48
1:A:1751:ARG:NE	1:A:1844:LEU:O	2.40	0.48
2:D:404:SER:O	2:D:405:ARG:CB	2.58	0.48
1:B:1748:GLY:N	1:B:1749:PRO:HD2	2.28	0.48
1:A:1752:ALA:HA	1:A:1845:TYR:CD2	2.49	0.48
1:C:1751:ARG:NE	1:C:1844:LEU:O	2.40	0.47
1:A:1700:THR:O	1:A:1703:TYR:HB3	2.14	0.47
1:C:1650:MET:N	1:C:1685:THR:HG21	2.29	0.47
1:B:1738:GLY:HA2	1:B:1746:HIS:CE1	2.49	0.47
1:C:1748:GLY:N	1:C:1749:PRO:HD2	2.29	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1672:HIS:O	1:C:1858:ILE:HG21	2.14	0.47
1:A:1822:HIS:CD2	1:B:1650:MET:HE1	2.50	0.47
2:E:402:GLU:C	2:E:403:TYR:CD2	2.87	0.47
1:B:1773:THR:CB	1:B:1774:ASN:OD1	2.63	0.46
1:C:1718:TRP:CZ2	1:C:1731:GLU:HG3	2.50	0.46
1:C:1682:GLU:CD	1:C:1682:GLU:H	2.19	0.46
1:C:1776:PRO:HG2	1:C:1778:ASP:OD1	2.15	0.46
1:A:1651:SER:OG	1:A:1684:THR:HA	2.14	0.46
1:B:1749:PRO:O	1:B:1753:ARG:HG3	2.15	0.46
1:C:1692:ASP:OD1	1:C:1696:VAL:N	2.42	0.46
1:A:1749:PRO:O	1:A:1753:ARG:HG3	2.15	0.46
1:C:1740:VAL:HG23	1:C:1741:VAL:HG23	1.97	0.46
1:A:1748:GLY:N	1:A:1749:PRO:HD2	2.31	0.46
1:B:1718:TRP:CZ2	1:B:1731:GLU:HG3	2.51	0.46
1:B:1775:MET:HG2	1:B:1776:PRO:O	2.16	0.45
1:C:1749:PRO:O	1:C:1753:ARG:HG3	2.17	0.45
1:A:1692:ASP:OD1	1:A:1696:VAL:N	2.43	0.45
2:E:401:GLY:HA2	2:E:402:GLU:HA	1.69	0.45
1:B:1740:VAL:HG23	1:B:1741:VAL:HG23	1.99	0.45
1:A:1738:GLY:HA2	1:A:1746:HIS:CE1	2.51	0.45
1:B:1774:ASN:N	1:B:1774:ASN:OD1	2.50	0.45
1:B:1751:ARG:NE	1:B:1844:LEU:O	2.40	0.45
1:A:1816:THR:O	1:A:1817:GLU:C	2.55	0.45
1:C:1776:PRO:HD2	1:C:1779:GLN:CG	2.47	0.44
2:E:402:GLU:O	2:E:403:TYR:CD2	2.70	0.44
1:A:1728:MET:HE3	1:C:1819:ASN:CB	2.47	0.44
1:C:1663:MET:N	1:C:1663:MET:SD	2.90	0.44
1:A:1773:THR:C	1:A:1775:MET:H	2.21	0.44
1:A:1656:GLY:N	2:D:406:SEP:O2P	2.40	0.44
1:A:1729:LEU:O	1:C:1819:ASN:CB	2.66	0.44
1:B:1816:THR:O	1:B:1817:GLU:C	2.56	0.44
1:A:1682:GLU:H	1:A:1682:GLU:CD	2.21	0.44
1:B:1651:SER:OG	1:B:1683:GLU:O	2.34	0.43
1:C:1776:PRO:HD2	1:C:1779:GLN:CB	2.48	0.43
1:A:1772:PHE:CD1	1:A:1777:THR:HG22	2.54	0.43
1:A:1731:GLU:OE1	1:C:1819:ASN:CB	2.67	0.43
1:C:1738:GLY:HA2	1:C:1746:HIS:CE1	2.53	0.43
1:A:1832:VAL:HG12	1:A:1855:ILE:HD12	2.01	0.43
1:A:1669:ALA:O	1:A:1670:ARG:C	2.57	0.43
1:B:1700:THR:O	1:B:1703:TYR:HB3	2.18	0.43
1:B:1663:MET:N	1:B:1663:MET:SD	2.92	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1718:TRP:CZ2	1:A:1731:GLU:HG3	2.53	0.43
1:C:1816:THR:O	1:C:1817:GLU:C	2.56	0.43
1:A:1663:MET:N	1:A:1663:MET:SD	2.92	0.42
1:C:1776:PRO:O	1:C:1779:GLN:N	2.53	0.42
1:C:1778:ASP:O	1:C:1779:GLN:C	2.55	0.42
1:C:1806:PRO:C	1:C:1807:ILE:HG13	2.40	0.42
1:B:1800:LEU:HD23	1:B:1800:LEU:N	2.34	0.42
2:E:402:GLU:HB3	2:E:403:TYR:CE1	2.54	0.42
1:C:1776:PRO:CB	1:C:1778:ASP:OD1	2.67	0.42
1:C:1700:THR:O	1:C:1703:TYR:HB3	2.20	0.42
1:C:1649:ARG:O	1:C:1650:MET:HE1	2.19	0.42
1:A:1771:PRO:HD2	1:A:1815:TRP:CD1	2.54	0.42
1:C:1769:TYR:O	1:C:1810:VAL:HG12	2.19	0.42
1:C:1648:LYS:HA	1:C:1649:ARG:HA	1.71	0.41
1:B:1773:THR:HB	1:B:1774:ASN:OD1	2.20	0.41
1:C:1651:SER:HA	1:C:1675:THR:HB	2.02	0.41
1:B:1752:ALA:HA	1:B:1845:TYR:CD2	2.56	0.41
1:A:1728:MET:HE2	1:C:1819:ASN:O	2.20	0.41
1:B:1759:LYS:HA	1:B:1845:TYR:OH	2.20	0.41
1:A:1646:VAL:HG23	1:A:1647:ASN:ND2	2.35	0.41
1:B:1769:TYR:O	1:B:1810:VAL:HG12	2.20	0.41
1:A:1769:TYR:O	1:A:1810:VAL:HG12	2.20	0.41
1:C:1649:ARG:O	1:C:1650:MET:SD	2.78	0.41
1:C:1826:GLN:CG	1:C:1859:PRO:HA	2.50	0.40
1:B:1832:VAL:HG12	1:B:1855:ILE:HD12	2.02	0.40
1:A:1698:GLU:OE2	2:D:408:THR:HG21	2.22	0.40
1:C:1832:VAL:HG12	1:C:1855:ILE:HD12	2.03	0.40
1:A:1821:PHE:CD1	1:A:1821:PHE:N	2.90	0.40
1:C:1759:LYS:HA	1:C:1845:TYR:OH	2.22	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1670:ARG:NH2	2:E:405:ARG:O[3_655]	2.18	0.02

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	212/214 (99%)	188 (89%)	24 (11%)	0	100 100
1	B	212/214 (99%)	187 (88%)	24 (11%)	1 (0%)	34 77
1	C	210/214 (98%)	186 (89%)	24 (11%)	0	100 100
2	D	6/11 (54%)	1 (17%)	2 (33%)	3 (50%)	0 0
2	E	8/11 (73%)	5 (62%)	3 (38%)	0	100 100
2	F	4/11 (36%)	1 (25%)	2 (50%)	1 (25%)	0 1
All	All	652/675 (97%)	568 (87%)	79 (12%)	5 (1%)	24 69

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	402	GLU
2	D	403	TYR
2	F	404	SER
2	D	405	ARG
1	B	1683	GLU

5.3.2 Protein sidechains

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	188/191 (98%)	178 (95%)	10 (5%)	28 67
1	B	189/191 (99%)	176 (93%)	13 (7%)	19 59
1	C	186/191 (97%)	175 (94%)	11 (6%)	24 65

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
2	D	6/8 (75%)	4 (67%)	2 (33%)	0 2
2	E	8/8 (100%)	6 (75%)	2 (25%)	1 4
2	F	5/8 (62%)	5 (100%)	0	100 100
All	All	582/597 (98%)	544 (94%)	38 (6%)	21 62

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1648	LYS
1	A	1649	ARG
1	A	1750	LYS
1	A	1774	ASN
1	A	1794	GLU
1	A	1799	THR
1	A	1821	PHE
1	A	1850	LEU
1	A	1852	THR
1	A	1854	LEU
1	B	1647	ASN
1	B	1649	ARG
1	B	1682	GLU
1	B	1750	LYS
1	B	1774	ASN
1	B	1777	THR
1	B	1794	GLU
1	B	1799	THR
1	B	1802	THR
1	B	1821	PHE
1	B	1850	LEU
1	B	1852	THR
1	B	1854	LEU
1	C	1648	LYS
1	C	1651	SER
1	C	1677	THR
1	C	1750	LYS
1	C	1794	GLU
1	C	1800	LEU
1	C	1802	THR
1	C	1821	PHE
1	C	1850	LEU
1	C	1852	THR

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Mol	Chain	Res	Type
1	C	1854	LEU
2	D	404	SER
2	D	405	ARG
2	E	402	GLU
2	E	403	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1647	ASN
1	A	1672	HIS
1	B	1672	HIS
1	C	1672	HIS
1	C	1848	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains (i)

3 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	SEP	D	406	2	7,9,10	0.77	0	8,12,14	2.50	4 (50%)
2	SEP	E	406	2	7,9,10	0.68	0	8,12,14	2.18	3 (37%)
2	SEP	F	406	2	7,9,10	0.62	0	8,12,14	1.65	2 (25%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SEP	D	406	2	-	0/5/8/10	0/0/0/0
2	SEP	E	406	2	-	0/5/8/10	0/0/0/0
2	SEP	F	406	2	-	0/5/8/10	0/0/0/0

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	D	406	SEP	OG-CB-CA	-4.46	104.38	108.26
2	E	406	SEP	OG-P-O1P	-3.44	98.43	107.08
2	D	406	SEP	OG-P-O1P	-3.13	99.23	107.08
2	D	406	SEP	O-C-CA	-2.91	117.93	125.72
2	F	406	SEP	O-C-CA	-2.69	118.50	125.72
2	F	406	SEP	OG-P-O1P	-2.51	100.79	107.08
2	E	406	SEP	O-C-CA	-2.38	119.33	125.72
2	D	406	SEP	O2P-P-O1P	2.45	118.61	110.63
2	E	406	SEP	O3P-P-O1P	3.53	122.16	110.63

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	406	SEP	1	0
2	E	406	SEP	3	0
2	F	406	SEP	1	0

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å^2)	Q<0.9
1	A	214/214 (100%)	0.11	0 100 100	94, 129, 177, 216	0
1	B	214/214 (100%)	0.45	13 (6%) 25 19	106, 167, 250, 286	0
1	C	212/214 (99%)	0.57	21 (9%) 9 9	134, 192, 227, 250	0
2	D	8/11 (72%)	-0.27	0 100 100	123, 165, 187, 214	0
2	E	10/11 (90%)	0.06	1 (10%) 9 9	164, 182, 206, 211	0
2	F	6/11 (54%)	-0.27	0 100 100	142, 162, 196, 212	0
All	All	664/675 (98%)	0.36	35 (5%) 30 23	94, 163, 232, 286	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	1716	TYR	5.9
1	C	1845	TYR	4.8
1	C	1715	SER	4.0
1	B	1840	ASP	4.0
1	B	1859	PRO	3.9
1	C	1712	TRP	3.8
1	B	1849	GLU	3.6
1	C	1689	MET	3.1
1	C	1709	GLY	3.1
1	C	1659	PRO	3.0
1	B	1707	ILE	3.0
1	B	1828	CYS	2.9
1	B	1700	THR	2.9
1	C	1676	LEU	2.8
1	B	1794	GLU	2.7
1	B	1703	TYR	2.6
1	C	1711	LYS	2.6
2	E	399	GLY	2.6
1	C	1660	GLU	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	1718	TRP	2.4
1	B	1795	LEU	2.4
1	C	1708	ALA	2.4
1	C	1657	LEU	2.4
1	C	1654	VAL	2.4
1	C	1667	LYS	2.2
1	B	1812	PRO	2.2
1	C	1690	LYS	2.2
1	C	1704	PHE	2.2
1	B	1807	ILE	2.2
1	B	1839	LEU	2.1
1	C	1656	GLY	2.1
1	C	1677	THR	2.1
1	B	1654	VAL	2.1
1	C	1759	LYS	2.1
1	C	1706	GLY	2.0

6.2 Non-standard residues in protein, DNA, RNA chains (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å^2)	Q<0.9
2	SEP	E	406	10/11	0.90	0.21	-	139,146,156,161	0
2	SEP	F	406	10/11	0.91	0.15	-	133,169,197,226	0
2	SEP	D	406	10/11	0.93	0.23	-	107,127,155,160	0

6.3 Carbohydrates (i)

There are no carbohydrates in this entry.

6.4 Ligands (i)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.